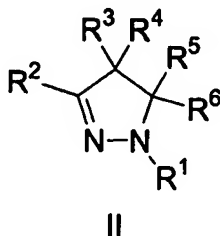


**In the claims:**

1.-2. (Previously cancelled)

3. (Currently amended) A compound of the Formula II,



wherein:

a is 0 or 1;

b is 0 or 1;

m is 0, 1, or 2;

R<sup>1</sup> is selected from:

- 1) (C=O)C<sub>1</sub>-C<sub>10</sub> alkyl,
- 7) (C=O)OC<sub>1</sub>-C<sub>10</sub> alkyl, and
- 8) (C=O)NR<sup>7</sup>R<sup>8</sup>,

said alkyl is optionally substituted with one or more substituents selected from R<sup>7</sup>; orR<sup>2</sup> is ~~selected from~~ phenyl;said phenyl is optionally substituted with one or more substituents selected from (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl, (C=O)<sub>a</sub>O<sub>b</sub>aryl, CO<sub>2</sub>H, halo, CN, ~~or (C=O)<sub>a</sub>NR<sup>9</sup>R<sup>10</sup>~~ or CHO;R<sup>3</sup> and R<sup>4</sup> are hydrogen;R<sup>5</sup> is selected from:

- 1) H, and
- 2) C<sub>1</sub>-C<sub>10</sub> alkyl,

said alkyl is optionally substituted with one or more substituents selected from R<sup>7</sup>;R<sup>6</sup> is phenyl:said phenyl is optionally substituted with one or more substituents selected from R<sup>7</sup>,

R<sup>7</sup> is:

- 1) (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 2) (C=O)<sub>a</sub>O<sub>b</sub>aryl,
- 3) CO<sub>2</sub>H,
- 4) halo,
- 5) CN,
- 6) OH,
- 7) O<sub>a</sub>(C=O)<sub>b</sub>NR<sup>9</sup>R<sup>10</sup>, and
- 8) CHO,

said alkyl, and aryl are optionally substituted with one, two or three substituents selected from R<sup>8</sup>;

R<sup>8</sup> is selected from:

- 1) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>1</sub>-C<sub>10</sub>)alkyl, wherein r and s are independently 0 or 1,
- 2) O<sub>r</sub>(C<sub>1</sub>-C<sub>3</sub>)perfluoroalkyl, wherein r is 0 or 1,
- 3) oxo,
- 4) OH,
- 5) halo,
- 6) CN,
- 7) (C<sub>2</sub>-C<sub>10</sub>)alkenyl,
- 8) (C<sub>2</sub>-C<sub>10</sub>)alkynyl,
- 9) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>3</sub>-C<sub>6</sub>)cycloalkyl,
- 10) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-aryl,
- 11) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-heterocyclyl,
- 12) (C=O)<sub>r</sub>O<sub>s</sub>(C<sub>0</sub>-C<sub>6</sub>)alkylene-N(R<sup>b</sup>)<sub>2</sub>,
- 13) C(O)R<sup>a</sup>,
- 14) (C<sub>0</sub>-C<sub>6</sub>)alkylene-CO<sub>2</sub>R<sup>a</sup>,
- 15) C(O)H,
- 16) (C<sub>0</sub>-C<sub>6</sub>)alkylene-CO<sub>2</sub>H, and
- 17) C(O)N(R<sup>b</sup>)<sub>2</sub>,
- 18) S(O)<sub>m</sub>R<sup>a</sup>, and
- 19) S(O)<sub>2</sub>NR<sup>9</sup>R<sup>10</sup>

said alkyl, alkenyl, alkynyl, cycloalkyl, aryl, and heterocyclyl is optionally substituted with up to three substituents selected from R<sup>b</sup>, OH, (C<sub>1</sub>-C<sub>6</sub>)alkoxy, halogen, CO<sub>2</sub>H, CN, O(C=O)C<sub>1</sub>-C<sub>6</sub> alkyl, oxo, and N(R<sup>b</sup>)<sub>2</sub>;

R<sup>9</sup> and R<sup>10</sup> are independently selected from:

- 1) H,
- 2) (C=O)O<sub>b</sub>C<sub>1</sub>-C<sub>10</sub> alkyl,
- 3) (C=O)O<sub>b</sub>C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 4) (C=O)O<sub>b</sub>aryl,
- 5) (C=O)O<sub>b</sub>heterocyclyl,
- 6) C<sub>1</sub>-C<sub>10</sub> alkyl,
- 7) aryl,
- 8) C<sub>2</sub>-C<sub>10</sub> alkenyl,
- 9) C<sub>2</sub>-C<sub>10</sub> alkynyl,
- 10) heterocyclyl,
- 11) C<sub>3</sub>-C<sub>8</sub> cycloalkyl,
- 12) SO<sub>2</sub>R<sup>a</sup>, and
- 13) (C=O)NR<sup>b</sup><sub>2</sub>,

said alkyl, cycloalkyl, aryl, heterocyclyl, alkenyl, and alkynyl is optionally substituted with one, two or three substituents selected from R<sup>8</sup>, or

R<sup>9</sup> and R<sup>10</sup> can be taken together with the nitrogen to which they are attached to form a monocyclic or bicyclic heterocycle with 5-7 members in each ring and optionally containing, in addition to the nitrogen, one or two additional heteroatoms selected from N, O and S, said monocyclic or bicyclic heterocycle optionally substituted with one, two or three substituents selected from R<sup>8</sup>;

R<sup>a</sup> is (C<sub>1</sub>-C<sub>6</sub>)alkyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, aryl, or heterocyclyl; and

R<sup>b</sup> is H, (C<sub>1</sub>-C<sub>6</sub>)alkyl, aryl, heterocyclyl, (C<sub>3</sub>-C<sub>6</sub>)cycloalkyl, (C=O)OC<sub>1</sub>-C<sub>6</sub> alkyl, (C=O)C<sub>1</sub>-C<sub>6</sub> alkyl or S(O)<sub>2</sub>R<sup>a</sup>.

4. (Currently amended) The compound according to Claim 3 or a pharmaceutically acceptable salt or stereoisomer thereof, wherein:

R<sup>1</sup> is selected from:

- 1) (C=O)C<sub>1</sub>-C<sub>10</sub> alkyl, and
- 4) (C=O)OC<sub>1</sub>-C<sub>10</sub> alkyl,

said alkyl, ~~aryl, cycloalkyl, and heterocyclyl~~ is optionally substituted with one, two or three substituents selected from R<sup>7</sup>;

R<sup>2</sup> is phenyl,

said phenyl is optionally substituted with one or more substituents selected from (C=O)<sub>a</sub>O<sub>b</sub>C<sub>1-C<sub>10</sub></sub> alkyl, (C=O)<sub>a</sub>O<sub>b</sub>aryl, CO<sub>2</sub>H, halo, CN,  $\Theta_a(C=O)_bNR^9R^{10}$  or CHO;

R<sup>3</sup> and R<sup>4</sup> are hydrogen;

R<sup>5</sup> is selected from:

- 1) H, and
- 2) C<sub>1</sub>-C<sub>10</sub> alkyl,

said alkyl is optionally substituted with one or more substituents selected from R<sup>7</sup>;

R<sup>6</sup> is phenyl:

said phenyl is optionally substituted with one or more substituents selected from R<sup>7</sup>, and R<sup>7</sup>, R<sup>8</sup>, R<sup>9</sup>, R<sup>10</sup>, R<sup>a</sup> and R<sup>b</sup> are as described in Claim 3.

5. (Previously cancelled)

6. (Previously amended) A compound selected from:

3-[1-acetyl-3-(2-chlorophenyl)-4,5-dihydro-1H-pyrazol-5-yl]phenol

3-[3-(2-chlorophenyl)-1-isobutyryl-4,5-dihydro-1H-pyrazol-5-yl]phenol

3-[1-acetyl-3-(2-chlorophenyl)-5-methyl-4,5-dihydro-1H-pyrazol-5-yl]phenol

3-[1-acetyl-3-(2,5-difluorophenyl)-4,5-dihydro-1H-pyrazol-5-yl]phenol

3-[1-Acetyl-3-(2-fluorophenyl)-4,5-dihydro-1H-pyrazol-5-yl]phenol

3-[1-Acetyl-3-(3-bromophenyl)-4,5-dihydro-1H-pyrazol-5-yl]phenol

3-[1-Acetyl-3-(2,3-dichlorophenyl)-4,5-dihydro-1H-pyrazol-5-yl]phenol

3-[1-Acetyl-3-(2,5-dichlorophenyl)-4,5-dihydro-1H-pyrazol-5-yl]phenol

3-[1-Propionyl-3-(2-chlorophenyl)-4,5-dihydro-1H-pyrazol-5-yl]phenol

3-[1-Isobutyryl-3-(2-chlorophenyl)-4,5-dihydro-1H-pyrazol-5-yl]phenol

1-Acetyl-3-(2-chlorophenyl)-5-phenyl-4,5-dihydro-1H-pyrazole

1-Acetyl-3-(3-chlorophenyl)-5-phenyl-4,5-dihydro-1H-pyrazole

1-Acetyl-3-(2,5-difluorophenyl)-5-phenyl-4,5-dihydro-1H-pyrazole

3-(2,5-difluorophenyl)- N,N-dimethyl-5-phenyl-4,5-dihydro-1H-pyrazole-1-carboxamide

3-(2,5-difluorophenyl)-5-(3-hydroxyphenyl)-N,N-dimethyl-4,5-dihydro-1H-pyrazole-1-carboxamide

3-(2,5-difluorophenyl)- N,N-diethyl-5-phenyl-4,5-dihydro-1H-pyrazole-1-carboxamide

3-(2,5-difluorophenyl)-5-phenyl-4,5-dihydro-1H-pyrazole-1-carboxamide

1-acetyl-3-(2,5-difluorophenyl)-5-methyl-5-phenyl-4,5-dihydro-1H-pyrazole

3-(2,5-difluorophenyl)-N,5-dimethyl-5-phenyl-4,5-dihydro-1H-pyrazole-1-carboxamide

3-(2,5-difluorophenyl)-N,N,5-trimethyl-5-phenyl-4,5-dihydro-1H-pyrazole-1-carboxamide

3-(2,5-difluorophenyl)-5-ethyl-N-methyl-5-phenyl-4,5-dihydro-1H-pyrazole-1-carboxamide

3-(2,5-difluorophenyl)-5-(hydroxymethyl)-N-methyl-5-phenyl-4,5-dihydro-1H-pyrazole-1-carboxamide

ethyl [3-(2,5-difluorophenyl)-5-methyl-5-phenyl-4,5-dihydro-1H-pyrazol-1-yl]acetate

ethyl [3-(2,5-difluorophenyl)-5-phenyl-4,5-dihydro-1H-pyrazol-1-yl]acetate

ethyl 2-[3-(2,5-difluorophenyl)-5-methyl-5-phenyl-4,5-dihydro-1H-pyrazol-1-yl]propanoate

3-(2,5-difluorophenyl)-5-[3-(dimethylamino)propyl]-N-ethyl-5-phenyl-4,5-dihydro-1H-pyrazole-1-carboxamide

3-(2,5-difluorophenyl)-N-ethyl-5-{3-[(1H-imidazol-2-ylcarbonyl)amino]propyl}-5-phenyl-4,5-dihydro-1H-pyrazole-1-carboxamide

5-(2-aminoethyl)-3-(2,5-difluorophenyl)-N-methyl-5-phenyl-4,5-dihydro-1H-pyrazole-1-carboxamide

5-(3-aminopropyl)-3-(2,5-difluorophenyl)-N-ethyl-5-phenyl-4,5-dihydro-1H-pyrazole-1-carboxamide

5-(3-aminobutyl)-3-(2,5-difluorophenyl)-N-ethyl-5-phenyl-4,5-dihydro-1H-pyrazole-1-carboxamide

5-[3-(benzoylamino)propyl]-3-(2,5-difluorophenyl)-N-ethyl-5-phenyl-4,5-dihydro-1H-pyrazole-1-carboxamide

3-(2,5-difluorophenyl)-5-[4-(dimethylamino)butyl]-N-ethyl-5-phenyl-4,5-dihydro-1H-pyrazole-1-carboxamide

3-(2,5-difluorophenyl)-5-[4-(dimethylnitro)but-1-yl]-N-ethyl-5-phenyl-4,5-dihydro-1H-pyrazole-1-carboxamide

5-[4-(benzylamino)butyl]-3-(2,5-difluorophenyl)-N-ethyl-5-phenyl-4,5-dihydro-1H-pyrazole-1-carboxamide

3-(2,5-difluorophenyl)-N-ethyl-5-phenyl-5-{4-[(pyridin-4-ylmethyl)amino]butyl}-4,5-dihydro-1H-pyrazole-1-carboxamide

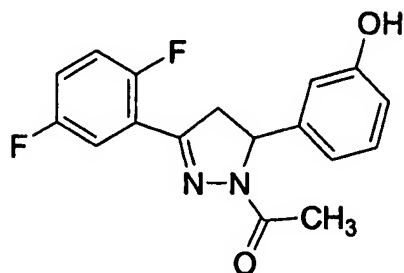
5-(3-amino-3-phenylpropyl)-3-(2,5-difluorophenyl)-N,N-dimethyl-5-phenyl-4,5-dihydro-1H-pyrazole-1-carboxamide

or a pharmaceutically acceptable salt or stereoisomer thereof.

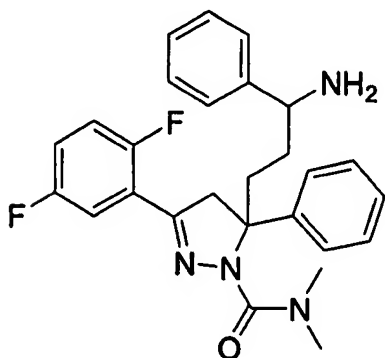
7. (Previously cancelled)

8. (Previously amended) The compound according to Claim 3 which is selected from:

3-[1-acetyl-3-(2,5-difluorophenyl)-4,5-dihydro-1H-pyrazol-5-yl]phenol

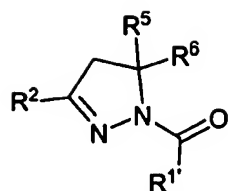


5-(3-amino-3-phenylpropyl)-3-(2,5-difluorophenyl)-N,N-dimethyl-5-phenyl-4,5-dihydro-1H-pyrazole-1-carboxamide

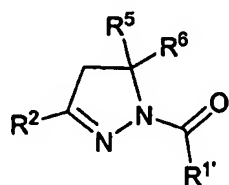


or a pharmaceutically acceptable salt or stereoisomer thereof.

9. (Previously amended) A compound selected from:



R <sup>2</sup>	R <sup>5</sup>	R <sup>6</sup>	R <sup>1'</sup>
2,5-dichlorophenyl	H	Ph	NMe <sub>2</sub>
2-fluoro-5-cyanophenyl	H	Ph	NMe <sub>2</sub>
2-fluoro-5-bromophenyl	H	Ph	NMe <sub>2</sub>
2-fluoro-5-chlorophenyl	H	Ph	NMe <sub>2</sub>
2-fluoro-5-nitrophenyl	H	Ph	NMe <sub>2</sub>



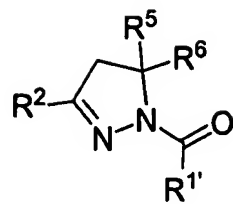
R <sup>2</sup>	R <sup>5</sup>	R <sup>6</sup>	R <sup>1'</sup>
2,5-difluorophenyl	H	3-hydroxyphenyl	NMe <sub>2</sub>
2,5-difluorophenyl	H	4-hydroxyphenyl	NMe <sub>2</sub>
2,5-difluorophenyl	H	3-aminophenyl	NMe <sub>2</sub>
2,5-difluorophenyl	H	3-(acetylamino)phenyl	NMe <sub>2</sub>



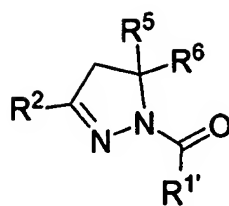
2,5-difluorophenyl

H

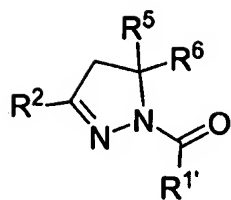
3-carboxyphenyl

NMe<sub>2</sub>

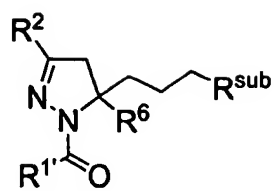
R <sup>2</sup>	R <sup>5</sup>	R <sup>6</sup>	R <sup>1'</sup>
2,5-difluorophenyl	H	Ph	
2,5-difluorophenyl	H	Ph	
2,5-difluorophenyl	H	Ph	



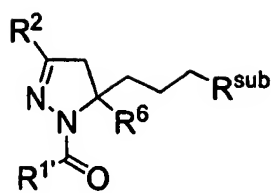
$R^2$	$R^5$	$R^6$	$R^{1'}$
2,5-difluorophenyl	H	Ph	
2,5-difluorophenyl	H	Ph	
2,5-difluorophenyl	H	Ph	



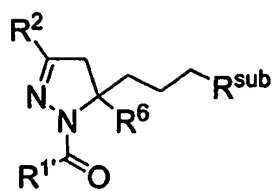
R <sup>2</sup>	R <sup>5</sup>	R <sup>6</sup>	R <sup>1'</sup>
2,5-difluorophenyl	H	Ph	
2,5-difluorophenyl	H	Ph	
2,5-difluorophenyl	H	Ph	
2,5-difluorophenyl	H	Ph	



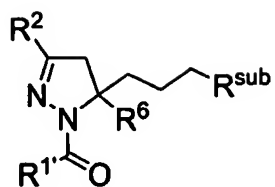
$R^2$	$R^{sub}$	$R^6$	$R^{1'}$
2,5-difluorophenyl	$NH_2$	Ph	
2,5-difluorophenyl	$NH_2$	Ph	
2,5-difluorophenyl	$NH_2$	Ph	



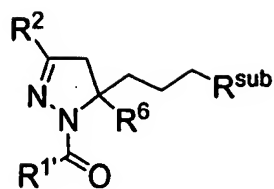
$R^2$	$R^{sub}$	$R^6$	$R^{1'}$
2,5-difluorophenyl	$NH_2$	Ph	
2,5-difluorophenyl	$NH_2$	Ph	
2,5-difluorophenyl	$NH_2$	Ph	
2,5-difluorophenyl	$NH_2$	Ph	



$R^2$	$R^{sub}$	$R^6$	$R^{1'}$
2,5-difluorophenyl	$NH_2$	Ph	
2,5-difluorophenyl	$NH_2$	Ph	
2,5-difluorophenyl	$NH_2$	Ph	

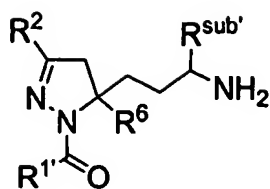


$R^2$	$R^{sub}$	$R^6$	$R^{1'}$
2,5-difluorophenyl	$NH_2$	3-hydroxyphenyl	$NMe_2$
2,5-difluorophenyl	$NH_2$	4-hydroxyphenyl	$NMe_2$
2,5-difluorophenyl	$NH_2$	3-aminophenyl	$NMe_2$
2,5-difluorophenyl	$NH_2$	3-(acetylamino)phenyl	$NMe_2$
2,5-difluorophenyl	$NH_2$	3-carboxyphenyl	$NMe_2$
2,5-difluorophenyl	$NH_2$	3-tetrazolylphenyl	$NMe_2$

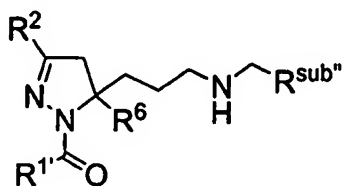


R <sup>2</sup>	R <sup>sub</sup>	R <sup>6</sup>	R <sup>1'</sup>
2,5-dichlorophenyl	NH <sub>2</sub>	Ph	NMe <sub>2</sub>
2-fluoro-5-cyanophenyl	NH <sub>2</sub>	Ph	NMe <sub>2</sub>
2-fluoro-5-bromophenyl	NH <sub>2</sub>	Ph	NMe <sub>2</sub>
2-fluoro-5-chlorophenyl	NH <sub>2</sub>	Ph	NMe <sub>2</sub>
2-fluoro-5-nitrophenyl	NH <sub>2</sub>	Ph	NMe <sub>2</sub>

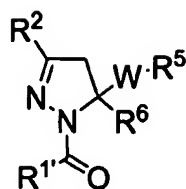




$R^2$	$R^{sub'}$	$R^6$	$R^{1'}$
2,5-difluorophenyl	phenyl	Ph	NMe <sub>2</sub>
2,5-difluorophenyl	4-trifluoromethylphenyl	Ph	NMe <sub>2</sub>
2,5-difluorophenyl	4-chlorophenyl	Ph	NMe <sub>2</sub>
2,5-difluorophenyl	CO <sub>2</sub> Me	Ph	NMe <sub>2</sub>
2,5-difluorophenyl	CONH <sub>2</sub>	Ph	NMe <sub>2</sub>



$R^2$	$R^{sub''}$	$R^6$	$R^{1'}$
2,5-difluorophenyl	phenyl	Ph	NMe <sub>2</sub>
2,5-difluorophenyl	4-trifluoromethylphenyl	Ph	NMe <sub>2</sub>
2,5-difluorophenyl	4-chlorophenyl	Ph	NMe <sub>2</sub>
2,5-difluorophenyl	CO <sub>2</sub> Me	Ph	NMe <sub>2</sub>
2,5-difluorophenyl	4-cyanophenyl	Ph	NMe <sub>2</sub>



$R^2$	$W-R^5$	$R^6$	$R^{1'}$
2,5-difluorophenyl	$-\text{CH}_2\text{CF}_2\text{CH}_2\text{NH}_2$	Ph	$\text{NMe}_2$
2,5-difluorophenyl	$-\text{CH}_2\text{OCH}_2\text{CH}_2\text{NH}_2$	Ph	$\text{NMe}_2$
2,5-difluorophenyl	$-\text{CH}_2\text{CH}_2\text{CH}(\text{CHF}_2)\text{NH}_2$	Ph	$\text{NMe}_2$
2,5-difluorophenyl	$-\text{CH}_2\text{OCF}_2\text{CH}_2\text{NH}_2$	Ph	$\text{NMe}_2$
2,5-difluorophenyl	$-\text{CH}_2\text{CH}_2\text{CF}_2\text{CH}_2\text{NH}_2$	Ph	$\text{NMe}_2$
2,5-difluorophenyl	$-\text{CH}_2\text{CH}_2\text{CH}_2\text{CH}(\text{CHF}_2)\text{NH}_2$	Ph	$\text{NMe}_2$
2,5-difluorophenyl	$-\text{CH}_2\text{CH}(\text{OH})\text{CH}_2\text{CH}_2\text{NH}_2$	Ph	$\text{NMe}_2$
2,5-difluorophenyl	$-\text{CH}_2\text{CH}(\text{OH})\text{CH}_2\text{NH}_2$	Ph	$\text{NMe}_2$
2,5-difluorophenyl	$-\text{CH}_2\text{C}(\text{O})\text{CH}_2\text{CH}_2\text{NH}_2$	Ph	$\text{NMe}_2$

or a pharmaceutically acceptable salt or stereoisomer thereof.

10. (Previously amended) A pharmaceutical composition that is comprised of a compound in accordance with Claim 3 and a pharmaceutically acceptable carrier.

11.- 36. (Previously cancelled)